

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
 Data File : BP026326.D
 Acq On : 15 Dec 2025 15:31
 Operator : RC/JU
 Sample : Q3839-19MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AM7MS

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 12/16/2025
 Supervised By :Jagrut Upadhyay 12/16/2025

Quant Time: Dec 15 15:57:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 15 14:15:35 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.649	152	213200	20.000	ng	0.00
21) Naphthalene-d8	10.413	136	807190	20.000	ng	0.00
39) Acenaphthene-d10	14.284	164	492568	20.000	ng	0.00
64) Phenanthrene-d10	17.084	188	974973	20.000	ng	0.00
76) Chrysene-d12	21.530	240	1085677	20.000	ng	0.00
86) Perylene-d12	24.824	264	1321757	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.278	112	615751	46.434	ng	0.00
7) Phenol-d6	6.837	99	467320	27.913	ng	0.00
23) Nitrobenzene-d5	8.790	82	977764	68.581	ng	0.00
42) 2,4,6-Tribromophenol	15.801	330	825909	126.169	ng	0.00
45) 2-Fluorobiphenyl	12.890	172	2539148	74.527	ng	0.00
79) Terphenyl-d14	19.825	244	4203399	77.139	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.208	88	82799	15.293	ng	95
3) Pyridine	3.602	79	173570	11.190	ng	97
4) n-Nitrosodimethylamine	3.508	42	95475	16.560	ng	# 94
6) Aniline	6.984	93	507674	23.146	ng	99
8) 2-Chlorophenol	7.225	128	487696	32.266	ng	98
9) Benzaldehyde	6.802	77	346378	38.689	ng	98
10) Phenol	6.867	94	212489	11.803	ng	98
11) bis(2-Chloroethyl)ether	7.078	93	520651	36.851	ng	98
12) 1,3-Dichlorobenzene	7.543	146	628968	38.357	ng	99
13) 1,4-Dichlorobenzene	7.678	146	634854	38.515	ng	98
14) 1,2-Dichlorobenzene	7.996	146	619567	39.152	ng	99
15) Benzyl Alcohol	7.890	79	293895	23.946	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.161	45	683354	34.603	ng	99
17) 2-Methylphenol	8.090	107	332132	27.022	ng	98
18) Hexachloroethane	8.708	117	224929	37.452	ng	97
19) n-Nitroso-di-n-propyla...	8.443	70	393968	37.885	ng	99
20) 3+4-Methylphenols	8.413	107	393602	23.140	ng	99
22) Acetophenone	8.461	105	829182	40.266	ng	# 98
24) Nitrobenzene	8.831	77	583439	40.453	ng	98
25) Isophorone	9.349	82	1130362	40.078	ng	99
26) 2-Nitrophenol	9.537	139	317197	43.123	ng	98
27) 2,4-Dimethylphenol	9.602	122	470786	36.325	ng	98
28) bis(2-Chloroethoxy)met...	9.831	93	690316	39.140	ng	100
29) 2,4-Dichlorophenol	10.066	162	531082	40.167	ng	99
30) 1,2,4-Trichlorobenzene	10.284	180	602134	44.415	ng	100
31) Naphthalene	10.466	128	1771275	40.333	ng	99
32) Benzoic acid	9.702	122	114380m	10.007	ng	
33) 4-Chloroaniline	10.572	127	425091	23.088	ng	99
34) Hexachlorobutadiene	10.749	225	383712	43.053	ng	99
35) Caprolactam	11.349	113	27612	5.875	ng	91
36) 4-Chloro-3-methylphenol	11.702	107	511710	36.560	ng	100
37) 2-Methylnaphthalene	12.078	142	1159103	37.268	ng	99
38) 1-Methylnaphthalene	12.302	142	1194221	39.352	ng	100
40) 1,2,4,5-Tetrachloroben...	12.454	216	705543	46.470	ng	99
41) Hexachlorocyclopentadiene	12.437	237	827305	84.448	ng	99
43) 2,4,6-Trichlorophenol	12.696	196	470401	45.484	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
 Data File : BP026326.D
 Acq On : 15 Dec 2025 15:31
 Operator : RC/JU
 Sample : Q3839-19MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AM7MS

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 12/16/2025
 Supervised By :Jagrut Upadhyay 12/16/2025

Quant Time: Dec 15 15:57:01 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 15 14:15:35 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.766	196	517975	44.836	ng	99
46) 1,1'-Biphenyl	13.101	154	1650374	44.131	ng	99
47) 2-Chloronaphthalene	13.143	162	1281520	43.631	ng	99
48) 2-Nitroaniline	13.349	65	361254	44.517	ng	99
49) Acenaphthylene	14.001	152	2202085	45.455	ng	99
50) Dimethylphthalate	13.737	163	1663084	44.708	ng	100
51) 2,6-Dinitrotoluene	13.854	165	371132	50.461	ng	99
52) Acenaphthene	14.348	154	1212018	42.812	ng	100
53) 3-Nitroaniline	14.190	138	243316	28.472	ng	97
54) 2,4-Dinitrophenol	14.401	184	378805	79.717	ng	97
55) Dibenzofuran	14.690	168	1957274	44.499	ng	99
56) 4-Nitrophenol	14.519	139	197517	26.228	ng	97
57) 2,4-Dinitrotoluene	14.660	165	519643	47.309	ng	94
58) Fluorene	15.348	166	1574772	45.100	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.925	232	495131	50.594	ng	97
60) Diethylphthalate	15.131	149	1709500	45.724	ng	99
61) 4-Chlorophenyl-phenyle...	15.354	204	790005	45.305	ng	96
62) 4-Nitroaniline	15.366	138	348540	39.224	ng	95
63) Azobenzene	15.654	77	1475119	47.387	ng	98
65) 4,6-Dinitro-2-methylph...	15.431	198	300929	47.164	ng	97
66) n-Nitrosodiphenylamine	15.572	169	1424145	47.138	ng	99
67) 4-Bromophenyl-phenylether	16.266	248	536949	49.216	ng	98
68) Hexachlorobenzene	16.384	284	633783	49.857	ng	97
69) Atrazine	16.548	200	560608	49.649	ng	98
70) Pentachlorophenol	16.742	266	881560	100.071	ng	99
71) Phenanthrene	17.125	178	2580477	46.729	ng	99
72) Anthracene	17.237	178	2632923	46.919	ng	100
73) Carbazole	17.501	167	2448066	46.238	ng	100
74) Di-n-butylphthalate	18.107	149	3045491	48.287	ng	99
75) Fluoranthene	19.225	202	3108066	46.476	ng	99
77) Benzidine	19.425	184	1249514	36.911	ng	99
78) Pyrene	19.601	202	3268688	45.100	ng	99
80) Butylbenzylphthalate	20.560	149	1478718	46.697	ng	97
81) Benzo(a)anthracene	21.507	228	3468163	46.335	ng	100
82) 3,3'-Dichlorobenzidine	21.436	252	966982	35.883	ng	99
83) Chrysene	21.577	228	3361028	47.559	ng	100
84) Bis(2-ethylhexyl)phtha...	21.454	149	2126595	44.937	ng	99
85) Di-n-octyl phthalate	22.724	149	3825314	46.497	ng	97
87) Indeno(1,2,3-cd)pyrene	28.559	276	4641606	46.604	ng	# 95
88) Benzo(b)fluoranthene	23.789	252	3743610	46.148	ng	99
89) Benzo(k)fluoranthene	23.854	252	3781372	46.398	ng	99
90) Benzo(a)pyrene	24.666	252	3653968	47.604	ng	99
91) Dibenzo(a,h)anthracene	28.642	278	3908453	47.988	ng	97
92) Benzo(g,h,i)perylene	29.712	276	3841882	47.447	ng	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

